10L081/voa

A-0U04-000071

#### ER PROGRAM DATA ASSESSMENT SUMMARY REPORT FORM

Bato	ch No. <u>8910L081</u>		Site Solar Ponds
Lab	oratory Roy F. Weston - L	ionville	No. of Samples/Matrix 4/Water
SOV	W # 10/86 (Rev. 2/88)	· · · · · · · · · · · · · · · · · · ·	Reviewer Org. TechLaw, Inc.
Sam	ple Numbers <u>SW094007,</u>	TB101089007, S	W095007, SW093007
	· · · · · · · · · · · · · · · · · · ·		
		Data As	ssessment Summary
		VOA	Comments
1.	Holding Times	A	Action Items 1,2; Comment 1
2.	GC/MS Tune/Instr. Perf.		
3.	Calibrations	A	Action Items 3,4; Comment 2
4.	Blanks	A	Action Item 5
5.	Surrogates	A	Action Items 6,7
6.	Matrix Spike/Dup.	X	Comment 3
7.	Other QC	V	
8.	Internal Standards	A	Action Items 6,7
9.	Compound Identification	X	Action Items 6,8; Comments 4,5
10.	System Performance	X	Comment 6
11.	Overa! Assessment	. <u> </u>	Data acceptable with qualifications.
	<ul> <li>V = Data had no problems.</li> <li>A = Data acceptable but qualified due to R = Data rejected.</li> <li>X = Problems, but do not affect data.</li> </ul>	o problems.	
Dot	• Auglity. Data contained in th	ic hatch were review	red and found to be acceptable with qualifications. Acceptable,
			impacted by the "Action Items" listed below are appropriately flagged.
	er to attached Results Summary Ta		The state of the s
			REVIEWS FIR CLASSIFICATION . 1
			By Langer (unin)
			Red 19/10/10

Action Items: 1) Non-detected results for aromatic compounds in samples SW093007, SW094007, and
TB101089007 are estimated and undetected (UJ) because holding times exceeded seven days.
2) All results in the dilution of sample SW095007 are rejected (R) because the holding time exceeded fourteen
days and the duration from VTSR to sample analysis exceeded ten days.
3) 4-Methyl-2-pentanone's RRF were less than 0.05 in all calibrations and 2-Butanone had either RRFs less
than 0.05, a %RSD or %D greater than 50% in the initial and continuing calibrations. As a result, the non-detected
results for these two compounds in all analyses are rejected (R).
4) Acetone and Methylene Chloride had %RSDs exceeding 50% in the 10/18/89 initial calibration and
Acetone's %D exceeded 50% in the 10/20/89 continuing calibration. Therefore, the non-detected results for
Methylene Chloride in sample SW094007 and Acetone in sample SW095007 are rejected (R). The positive results
for Acetone in the other three samples are estimated (J). Positive results for Methylene Chloride in samples
SW093007 and TB101089007 would be estimated (J) had blank criteria been met. See Action Item 5. The positive
result for Methylene Chloride in sample SW095007 would be estimated however, all results are rejected (R) for this
sample. See Action Item 6.
5) As a result of method blank contamination, the positive results for Methylene Chloride in samples
SW093007 and TB101089007 are estimated and undetected (UJ) according to the Functional Guidelines criteria
(5x and 10x rules).
6) Surrogate recoveries exceeded 300% and all three internal standard areas exceeded criteria in sample
SW095007. In addition, unknown peaks were evident in the chromatogram, including one enormous peak, all of
which were not quantitated or reported. Furthermore, the quantitation report had positive results for several
compounds that were changed to non-detects without being initialed or dated. None of these peaks or compounds,
including Methylene Chloride which exceeded calibration range, were evident in the dilution. As a result of these
contributory problems, all results in sample SW095007 are rejected (R).
7) All three internal standard areas exceeded criteria in samples SW093007, SW094007, and TB101089007
and sample SW094007 also had surrogate recoveries outside criteria. Therefore, all results in these three samples
are estimated (J) or estimated and undetected (UJ).

Action Items: (cont) 8) The positive result for Chloroform in sample SV	W094007 is estimated (J) because
mass spectral data was not submitted to confirm its identity. The value is report	ted because of a good RRT match
with the 12 hour standard.	
Comments: 1) Sample reanalyses were conducted beyond the fourteen day	and ten day contract holding times.
Therefore, the original analyses data are reported on the Data Summary Table.	
2) Other compounds whose %RSD or %D exceeded criteria in the calibrat	ions were undetected in the samples
and no action is necessary.	
3) No MS/MSD analysis was performed with these samples to determine i	f possible matrix effects contributed
to surrogate and internal standard problems. No further action is necessary as a	Il sample results are qualified.
4) The reference spectrum for Acetone did not match NBS standard spectr	um.
5) Vinyl Acetate's identity in sample SW095007 is not confirmed by the n	nass spectrum provided. However
all results in this sample are rejected.	
6) The surrogate and internal standard area difficulties may indicate a serio	ous system problem.
Note: Data Summary Tables are attached.	
William T Fea	6/25/80
Reviewer Signature	Date

Page 1 of 1

Low Water TABLE #: 8910L081
SITE NAME: Solar Ponds
CLP VOLATILE ORGANIC ANALYSIS:

ANALYTICAL RESULTS (ppb)

Semple Name         CRACK         SINCROSOTY         TRITOROGEON         SINCROSOTY	Sample Location									
mythle Diete         Method Blank         101/1089         101/1089         101/1089         101/1089         101/1089         101/1089         101/1089         101/1089         101/1089         101/1089         101/1089         101/1089         101/1089         101/1089         201/1089	Sample Number	VBLK044	SW083007	SW094007	TB101089007	SW095007	VBLK058	SW095007DL		
Participation   Participati	Sampling Date		10/10/89		10/10/89	10/10/89		10/10/89		
Maile   Mail		$\vdash$			Trip Blank			2x Dilution		
Control Registrate   Control						2	2	8		
The contraction   The contra			<b>₩</b> 6+	3	3	>	3	=		
The control of the	Bromomethane	10	10 W A	3	3	,		-		
No. of the contraction   10   M   10	Vinyl chloride	10	4 W 0t	3	3			5		
Charlesteeled   Charlesteele	Chloroethane	10		3	3	5		þ		
State   Stat	Methylene chloride	9	3	Þ	3			5		
December   S	Acetone	10	٦		4 C 7	5		5		
Obsidercontribute         S         UA         S         UA         S         UA         S         UA         C         UA <th< th=""><th>Carbon disulfide</th><th>9</th><th></th><th>3</th><th>8 W A</th><th>5</th><th></th><th>5</th><th></th><th></th></th<>	Carbon disulfide	9		3	8 W A	5		5		
Ubblishocontriane         S         UM A         S         UM A         S         UM A         S         U B         T <th< th=""><th>1,1-Dichloroethene</th><th>5</th><th></th><th>3</th><th>8 W A</th><th>5</th><th></th><th>5</th><th></th><th></th></th<>	1,1-Dichloroethene	5		3	8 W A	5		5		
Controlled   Control   C	1,1-Dichloroethane	2	3	3	3	1		5		
Decidential	1,2-Dichloroethene (Total)	5	3	3	3	1				
Contractive   Contraction	Chloroform	5	3	5	3	1		b		
10	1,2-Dichloroethane	2		3		1		5		
1.7     1.2	2-Butanone	10	•	b		1		ŀ		
Obout letrachloride         5         W A         6         M A <th>1,1,1-Trichloroethane</th> <th></th> <th>3</th> <th>3</th> <th>3</th> <th>j</th> <th></th> <th>1</th> <th></th> <th></th>	1,1,1-Trichloroethane		3	3	3	j		1		
by accelate         10         HO         10         HO         10         HO         2         R         20         R         10           Amodiciticommentarie         5         UM A         5         UM A         5         UM A         5         U A         10         10           Amodiciticommentarie         5         UM A         5         UM A         5         UM A         5         U B         10           4.3. Dichloroproparie         5         UM A         5         UM A         5         UM A         5         U B         10           Amoroditionmenthane         5         UM A         5         UM A         5         UM A         5         UM A         10         UM A         10         UM A         10         UM A         10         UM A         5         UM A	Carbon tetrachloride	2	3	3	3	1				
Decoderation   Deco	Vinyl acetate	10		3	3	Į				
10   10   10   10   10   10   10   10	Bromodichloromethane	5		3	3	1		1		
1,3.Dichkiocopropene	1,2-Dichloropropane	5		3	3	b		10 U R		
Connoctioned personal processes of personal processes and personal per	cis-1,3-Dichloropropene	5	3	3	3	5		-		
xomodloromethane         5         W         A         5         W         A         5         W         A         5         W         B         W         B         W         B         W         B         W         B         W         B         W         B         W         B         W         B         W         B         W         B         W         B         W         B         W         B         W         B         W         B         W         B         B         W         B         W         B	Trichloroethene	2	3	3	3			5		
2-Trichkoroethane         5         W         5         W         6         W         6         W         6         W         6         W         6         W         6         W         6         W         6         W         7         W         7         W         7         W         7         10         W         7         W         7         W         7         10         W         7         W         7         10         W         <	Dibromochloromethane	5	, ,	3	3	I		5		
Receive the contraction of the contract	1,1,2-Trichloroethane	2		3	3	1		ı		
re-1,3-Dichloropcropene         5         W A	Вепzеле	5	,	3	3	J				
Actify12-pentlanone         5 W A <th>trans-1,3-Dichloropropene</th> <th>2</th> <th>3</th> <th>3</th> <th>3</th> <th>1</th> <th></th> <th>5</th> <th></th> <th></th>	trans-1,3-Dichloropropene	2	3	3	3	1		5		
Adativit-2-pentianone         10         NR         10         UR         5         UR         7         10           Aybenzene         5         UR         5         UR         5         UR         5         UR         5         UR         5         UR         10         T         10         T         T         T         T         T         T         T         T         T         T         T         T         T         T         T         T         <	Bromoform	2	3	3	3	1		ı		
texanone         10         M A         5         M A	4-Methyl-2-pentanone	10		כ	כ			ĺ		
Leach Incomplementation of three Parameters         5 W A </th <th>2-Hexanone</th> <th>10</th> <th></th> <th>3</th> <th>3</th> <th></th> <th></th> <th>1</th> <th></th> <th></th>	2-Hexanone	10		3	3			1		
2.2-Tertrachlorocethane         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         A         A         B         B         A         A         B	Tetrachloroethene	5		3	3			ב		
Unene         5         W         A         5         W         A         5         W         A         5         W         B         T <th>1,1,2,2-Tetrachloroethane</th> <th>5</th> <th>3</th> <th>3</th> <th>3</th> <th></th> <th></th> <th>5</th> <th></th> <th></th>	1,1,2,2-Tetrachloroethane	5	3	3	3			5		
borobenzene         5         W         A         5         W         A         5         W         B         F         W         B         F         W         B         F         W         B         F         W         B         F         W         B         F         W         B         F         W         B         F         W         B         F         W         B         F         W         B         F         W         B         F         W         B         F         W         B         F         F         F         B         F <th< th=""><th>Toluene</th><th>5</th><th>3</th><th>3</th><th>3</th><th>Ι.</th><th></th><th>ı</th><th></th><th></th></th<>	Toluene	5	3	3	3	Ι.		ı		
ybenzene         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         5         W         A         10           ind Organic         ind Organic         ind Organic         E         24         8         7         247         3         0           Indicates the compound was not detected above the Required Outentified during the quality control review.         8         7         247         3         0           Cuantitation is approximate due to limitations identified during the quality control review.         Exceeds calibration in the programs per Liter (ug/L), Parts per billion (ppb).         7         247         X	Chlorobenzene	5		3	3			10 U R		
renee (Total)         5 W A         5 W A         5 W A         5 W A         5 W A         5 W A         10           Ital Organic         Ital Organic         6         24         8         7         247         9           Indicates the compound was not detected above the Required Oueritikation Limit.         6         24         8         7         247         9           Outsuitiation is approximate due to limitations identified during the quality control review.         Exceeds calibration range, dilute & reanalyze.         V           Exceeds calibration Limit in Micrograms per Liter (ug/L), Parts per billion (ptb).         R         A	Ethylbenzene	5		3	3	Ι.		0 10 ∪		
Secondary   Seco	Styrene	5		3	3	J.,		5		
ncentration (pbb)  Countration (pbb)  Exceeds calibration range, dilute & reanalyze.  Exceeds calibration Limit in Micrograms per Liter (ug/L), Parts per billion (pbb).	Xylenes (Total)	2	3	3	3	b		5		
noentration (ppb)  Indicates the compound was not detected above the Required Quantitation Limit.  Quantitation is approximate due to limitations identified during the quality control review.  Exceeds calibration range, dilute & reanalyze.  Exceeds calibration Limit in Micrograms per Liter (ug/L), Parts per billion (ppb).	Total Organic									
Indicates the compound was not detected above the Required Quantitation Limit.  Quantitation is approximate due to limitations identified during the quality control review.  Exceeds calibration range, dilute & reamalyze.  A  (QL. Contract Required Quantitation Limit in Micrograms per Liter (ug/L), Parts per billion (ppb).		9	24	1	7	247		0		-
> <b>«</b> α		setected above the Req	luired Quantitation L	lmit.				DQ Data Qualifier	alifier	
< α	J Cuantitation is approximate due to	limitations identified du		rol review.						
Œ	E Exceeds calbration range, dilute &	k reanalyze.							Acceptable with qualifications	
	CROL. Contract Required Quantitation	n Limit in Micrograms p	ver Litter (ug/L), Parts	s per billion (ppb).				R Rejected		

L081L/temp10

J Quantitation is approximate due to limitations identified during the quality control review.

E. Exceeds calibration range, dilute & reanalyze.

CROL. Contract Required Quantitation Limit in Micrograms per Liter (ug/L), Parts per billion (ppb).

### ER DEPARTMENT DATA ASSESSMENT SUMMARY REPORT FORM

Bato	ch No. 8910L081		Site Solar Ponds	
Lab	oratory Roy F. Weston - Li	onville	No. of Samples/Matrix <u>3/V</u>	Vater
SOV	<b>V</b> # 10/86 (Rev. 2/88)		Reviewer Org. <u>TechLaw, I</u>	nc.
Sam	ple Numbers <u>SW094007, S</u>	W095007, SW09	93007	
			·	
		Data As	sessment Summary	
		BNA	Comments	
1.	Holding Times			
2.	GC/MS Tune/Instr. Perf.			
3.	Calibrations	A	Action Item 1; Comment 1	
4.	Blanks	A	Action Item 2; Comment 2	
5.	Surrogates	X	Comment 3	
6.	Matrix Spike/Dup.	X	Comment 4	
7.	Other QC			e de la composição de l
8.	Internal Standards	<u> </u>		
9.	Compound Identification	X	Comment 5	
10.	System Performance	Y		
11.	Overall Assessment	A	Data Acceptable with Qualifications	
	V = Data had no problems.  A = Data acceptable but qualified due to  R = Data rejected.	problems.		
	X = Problems, but do not affect data.			
			ed and found to be acceptable with qualification	
	may be used provided that individually to attached Data Summary Table		by the "Action Items" listed below are appropria	uciy maggeo.
71/61	or to areaction water summary ravis	/3.j		<del></del>

Action Items: 1) The %RSD for Benzyl Alcohol exceeded 50% in the initial	calibration. The non-detected
Benzyl Alcohol results are rejected (R) in all samples.	
2) As a result of method blank contamination, the positive results for Di-n-but	ylphthalate and
Bis(2-ethylhexyl)phthalate in all samples is estimated and undetected (UI) accor-	ding to the Functional Guidelines
criteria.	
Comments: 1) Several compounds exceeded the %RSD and %D criteria in the	e 10/17/89 initial and the 10/19/89
continuing calibrations. Furthermore, Benzoic Acid's RRF was less than 0.05 in	this continuing calibration. The
only sample affected is the method blank and, therefore, data is not qualified.	
2) The method blank was analyzed on a different instrument than the samples.	Contamination found
in the blank only relates to possible contaminants in the samples introduced in the	e extraction process and not the
analysis itself.	
3) Two surrogates fell outside the % recovery criteria for the blank matrix spil	ke duplicate sample analysis.
4) A blank should not be used for the MS/MSD analysis since it is not represe	ntative of the sample matrix.
Moreover, six compounds exceeded the % recovery in the BS and BSD samples.	. No action is necessary
because results are not qualified based on MS/MSD data alone.	
5) Samples SW093007 and SW095007 contained four TICs and sample SW09	94007 contained five TICs.
Note: Data Summary Tables are attached.	
Lisa Contress - Herdler Reviewer Signature	7-//-90 Date
Reviewer Signature	Date

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Solar Ponds SITE NAME: Solar Pon CLP SEMIVOLATILE ANALYSIS: TABLE #:

8910L081

Low Water

ANALYTICAL RESULTS (ug/L)

Val)   CROL   DO   DO   DO   O   O   O   O   O   O	Sample Location								
Method Blantk   1071089	Sample Number		SBLK 967	77	SW095007	SW093007			
CFRQL   Method Blank   Method Blan	Sampling Date				10/10/89	10/10/89			
CFOL	Remarks		Method Blank						
Degree   Log   Degree   Degr	Semivolatiles (BNA)	CROL							
10	Organic Compound	ug/L	DO		8				
10	Phenol	10		ם	þ	10 U			
10	Bis(2 - Chloroethyl)ether	10		5	5	5			
10	2 - Chlorophenol	10		5	5	5			
10   10   V   10	1,3 - Dichlorobenzene	10		5	5	ł			
10	1,4 - Dichlorobenzene	10		5	>	ı			
10	Benzył alcohol	10		5	5	ı			
10   10   10   10   10   10   10   10	1,2 - Dichlorobenzene	10		5	5	ı			
10   10   10   10   10   10   10   10	2 - Methylphenol	10		5	5	5			
10	3is(2-chloroisopropyf)ether	10		5	b	1 .			
amine 10 10 0 V	t - Methylphenol	10		b	5	1			
10	V-Nitroso-di-n-propylamine	10		ລ	ם	1			
10	lexachloroethane	10		כ	ח				
10	litrobenzene	10		Ы	1	כ			
10	saphorone	9		כ		ח			
10   10   V   V   10   V   10   V   10   V   V   V   V   V   V   V   V   V	- Nitrophenol	10		D	1	D			
So U V   So U V   So U V	,4 - Dimethylphenol	9		5					
the finance 10 10 1 V 10 U V 1	enzoic Acid	SS		_	ם				
nee         10         V         10         V         10         V           stee         10         V         10         V         10         V           red         10         V         10         V         10         V           red         10         V         10         V         10         V           red         10         V         10         V         10         V           adiene         50         V         50         V         50         V           adiene         50         V         50         V         50         V           adiene         50         V         50         V         0         V           adien         50         V         50         V         V <t< td=""><td>is(2-Chloroethoxy)methane</td><td>9</td><td></td><td>_</td><td>b</td><td>ח</td><td></td><td></td><td></td></t<>	is(2-Chloroethoxy)methane	9		_	b	ח			
nne         10         V         10         V         10         V           10         10         V         10         V         10         V           20         10         V         10         V         10         V           30         10         V         10         V         10         V           4         10         V         10         V         10         V           50         10         V         10         V         10         V           10         10         V         10         V         10         V           10         10         V         10         V         V         V           10         10         V         10         V         V         V           10         V         10         V         10         V         V <t< td=""><td>,4 - Dichlorophenol</td><td>10</td><td></td><td>5</td><td>5</td><td>2</td><td></td><td></td><td></td></t<>	,4 - Dichlorophenol	10		5	5	2			
10   10   10   10   10   10   10   10	2,4 - Trichlorobenzene	9		٥	Ы	٥			
10   10   10   10   10   10   10   10	aphthalene	10		Ы	_	כ			
Adiene 10 10 10 10 10 10 10 10 10 10 10 10 10	- Chloroaniline	10		D	Ы	ם			
Adiene 10 10 10 V 10 U	lexachlorobutadiene	9		5	Ы	٥			
Adiene 10 10 0 V 10 U V	-Chloro-3-methylphenol	<b>£</b>		ם	Э	٥			
Adiene 10   10 U V	-Methylnaphthalene	9		Ы	5	5			
10   10   10   10   10   10   10   10	exachlorocyclopentadiene	ç		5	5	5			
50   0   50   0   0   0   0   0   0	,4,6-Trichlorophenol	9		٦	5	٥			
10   10   V   10	,4,5-Trichlorophenol	ន		٦	اد	b			
50   10   50   10   10   10   10   10	-Chloronaphthalene	0		5	5	5			
10   10   10   10   10   10   10   10	-Nitroaniline	50		ם	ח	כ			
10   10   10   10   10   10   10   10	Jimethyl phthalate	10		ם	n	ם			
10   10   10   10   V	\cenaphthylene	9		n	b	ם			
50 U V 50 U V DQ DQ DQ Station range.  So the compound was not detected above the Instrument Quantitation Limit.  So the compound was not detected above the Instrument Quantitation Limit.  V V V Detection is exproximate due to limitations identified during the quality control review.	,6-Dinitratoluene	9		_	٥	Ы			
8 > ∢ a	-Nitroaniline	20		Ы	כ	Ы			
> < a		٠					ualitier		
<b>&lt; a</b>		ot detected a	above the instrumer	nt Quantitation Limit					
Q	Ouantitation is approximate dux	e to limitation	s identified during	the quality control re	eview.		able with qualification	92	
<b>.</b>	CROL Contract Required Quantita	tion Limit in N	Vicrograms per Lite	r (ug/L), Parts per E	illion (ppb).				

U indicates the compound was not detected above the instrument Quantitation Limit.

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Solar Ponds 8910L081 SITE NAME:

TABLE #:

CLP SEMIVOLATILE ANALYSIS:

Low Water

ANALYTICAL RESULTS (ug/L)

Data Qualifier > 0 20 0 30 0 30 5 0 0 0 10 U V 20 U V 10 U V 10 U V 10 W A 10 U V 10 U V 10 U V 15 W A 10 U V 10 U V 10 U 10 O 10 C 10 U 10 U 10 U 우 우 SW093007 10/10/89 0 g 10 U V 10 U V ∪ 01 20 U V 10. U 10 U A V 10 U V 5 5 5 5 5 5 1 J A 10 U V 2 2 2 ⊃ 8 50 U 10 C 10 U 10 U SW095007 10/10/89 2 C C 2 S 2 U 2 V 10 U V 14 W A > 0 20 20 10 U V 10 U 10 U V 10 W A 10 U V 10 U 10 10 10 10 U 9 0 10 C ⊃ 83 2 28 10 U 10 U 10 U 10 U 10 U SW094007 10/10/89 0 8 å t ppb Method Blank 4 ppb SBLK 967 10 10 우우 5 5 2 5 우 ಜ જ 9 우 우 우 10 ន 9 9 9 10 10 2 Chlorophenyl-phenyl ether 4-Bromophenyl phenyl ether ,6-Dinitro-2-methylphenol Bis(2-ethylhexyl)phthalate V-Nitrosodiphenylamine Benzo(a)pyrene Indeno(1,2,3-cd)pyrene Dibenz(a,h)anthracene 3,3'-Dichlorobenzidine Total semivolatile BNA Benzo(b)fluoranthene Butylbenzyl phthalate Benzo(k)fluoranthene Hexachlorobenzene Benzo(a)anthracene Benzo(g,h,i)perylene Semivolatiles (BNA) Organic Compound Pentachlorophenol Di-n-butyl phthalate Di-n-octyl phthalate 2,4-Dinitrotoluene concentration (ppb) 2,4-Dinitrophenol Sample Location Diethyl phithalate Sample Number Sampling Date Acenaphthene -Nitrophenol Phenanthrene Dibenzofuran 4-Nitroaniline Fluoranthene Anthracene Remarks Fluorene Chrysene Pyrene

OL081.WK1

Acceptable with qualifications Rejected

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Exceeds calibration range.

Indicates the compound was not detected above the Instrument Quantitation Limit.

CROL. Contract Required Quantitation Limit in Micrograms per Liter (ug/L), Parts per Billion (ppb). Quantitation is approximate due to limitations identified during the quality control review.

# ER DEPARTMENT DATA ASSESSMENT SUMMARY REPORT FORM

Bato	ch No. <u>8910L081</u>			Site Solar Ponds	
Lab	oratory <u>Roy F. Weston - I</u>	Lionville		No. of Samples/Matrix _	3/Water
SOV	W # 10/86 (Rev. 2/88)			Reviewer Org. TechLaw	Inc.
Sam	ple Numbers <u>SW094007</u> .	SW095007, SW09	93007		
		Data Ass	essment	Summary	
		Pesticides/PCB		Comments	
1.	Holding Times	<u>V</u>		· · · · · · · · · · · · · · · · · · ·	
2.	Instrument Performance	X	Comme	ent 1	
3.	Calibrations	X	Commo	ent 2	<u> </u>
4.	Blanks	V			
5.	Surrogates	X	Commo	ent 3	
6.	Matrix Spike/Dup.	X	Comme	ent 4	· · · · · · · · · · · · · · · · · · ·
7.	Other QC	X	Comme	ent 5	
8.	Compound Identification	X	Commo	ent 6	<u> </u>
9.	System Performance	<u>v</u>		And the second s	
10.	Overall Assessment	<u> </u>	Data is	valid.	
	<ul> <li>V = Data had no problems.</li> <li>A = Data acceptable but qualified due t</li> <li>R = Data rejected.</li> <li>X = Problems, but do not affect data.</li> </ul>	o problems.			
Dat	a Quality: <u>Data contained in th</u>	nis batch were reviewed	d and foun	d to be valid. (Refer to attached I	Data Summary Table.)

Comments: 1) In the continuing calibration of 10/26/89 (0419) three compounds were outside of their retention
time windows. In the continuing calibration of 10/26/89 (0452) one compound was outside of its retention time
window. These calibrations were on the confirmation column. Expanded windows were used to evaluate the
chromatograms and no action is necessary.
2) The calibration factors of the multiresponse compounds (Aroclors and Toxaphene) are not reproducible. It
is suspected that multiple peaks were used to calculate the calibration factors. There were no positive results for
these compounds and no action is necessary.
3) Surrogate Recoveries were high in the method blank PBLK966 and sample SW093007. No action is
necessary.
4) No matrix spike/matrix spike duplicate data was provided with this batch. Although blank spike data was
included, this does not represent a sample matrix. In this blank spike data, Lindane's (gamma-BHC) RPD and
Percent Recoveries in both the BS/BSD exceeded the criteria. No action is necessary because results are not
qualified due to MS/MSD alone.
5) Although 990 milliliters of sample were extracted, the CRQLs were not adjusted accordingly. Therefore,
the CROLs were recalculated and the changed results reported on the Data Summary Table.
6) The positive result for DDT in the method blank is questioned. The retention time of DDT on the
confirmation column shifted, which is not a problem; however, it shifted early rather than late like the surrogate. In
addition the quantities generated from the two columns were very different. Nevertheless, the result is well below
the CRQL and no action is taken.
Note: Data Summary Tables are attached.
Reviewer Signature 7-10-90  Date
Keviewer Signature Date

Solar Ponds TABLE #: 8 SITE NAME:

8910L081

CLP PESTICIDES/PCB ANALYSIS: Low Water

Page 1 of 1

ANALYTICAL RESULTS (ug/L)

- 1 - 1 - 1 - 2								
Sample Location			1					
Sample Number		PBLK966	SW094007	SW095007	SW093007			
Sample Date			10/10/89	10/10/89	10/10/89			
Remarks		Method Blank						
Chlorinated Pesticides	CROL							
Analyte	ng√L	Ø	Ø	8	8			-
alpha - BHC	0.050		0.051 U V	0.051 U V	0.051 U V			
beta - BHC	0.050		0.051 U V	0.051 U V	0.051 U V			
detta - BHC	0.050		0.051 U V	0.051 U V	0.051 U V			
gamma - BHC (Lindane)	0.050		0.051 U V	0.051 U V	0.051 U V			
Heptachlor	0.050		ວ	0.051 U V	0.051 U V			
Aldrin	0.050		0.051 U V	0.051 U V	0.051 U V			
Heptachlor epoxide	0.050		0.051 U V	0.051 U V	0.051 U V			
Endosulfan I	0.050		0.051 U V	0.051 U V	0.051 U V			
Dieldrin	0.10		0.10 U V	0.10 U V	0.10 V V			
4,4' - DDE	0.10		0.10 U V	0.10 U V	0.10 V			
Endrin	0.10		0.10 U V	0.10 U V	0.10 V			
Endosulfan II	0.10		0.10 U V	0.10 U V	0.10 V			
4,4' - DDD	0.10	979 ddd 6.70	0.10 U V	0.10 U V	0.10 V			
Endosultan sultate	0.10		0.10 U V	0.10 U V	0.10 U V			
4,4' - DDT	0.10	0.010 ppb	0.10 U V	0.10 U V	0.10 U V			
Methoxychlor	0.50		5	0.51 U V	0.51 U V			
Endrin Ketone	0.10		Þ	0.10 U V	0.10 U V			
aipha - Chlordane	0.50		>	0.51 U V	0.51 U V			
gamma - Chlordane	0.50		0.51 U V	0.51 U V	0.51 U V			
Toxaphene	1.0		1.0 V	1.0 U V	1.0 V			
Aroclor - 1016	0.50		0.51 U V	0.51 U V	0.51 U V			
Arodor - 1221	0.50		ר	0.51 U V	0.51 U V			
Arodor - 1232	0.50		b	0.51 U V	0.51 U V			
Arodor - 1242	0.50	,	ם	0.51 U V	0.51 U V			
Arodor - 1248	0.50		0.51 U V	0.51 U V	0.51 U V			
Arodor - 1254	1.0		1.0 U V	1.0 U V	1.0 U V			
Arodor - 1260	1.0		1.0 U V	1.0 U V	1.0 U V			
Total Chlorinated Pesticides (ppb)		08.0	0	0	0			
E Exceeds calibration range.						DO Data Qualifier		

Indicates the compound was not detected above the Instrument Quantitation Limit

J Quantitation is approximate due to limitations identified during the quality control review CRQL. Contract Required Detection Limit in Micrograms per Liter (ug/L), Parts per Billion (ppb)

10L081/pcb

A Acceptable with qualifications Rejected

404161089001 Her	WESTON Analytics Use Only Samples Were:	Notes:	NOTES: 3 Received Broken	Sealed)  Y  NOTES:	A Property Preserved  V N N N N N N N N N N N N N N N N N N	5 Received Within Thoding Times Y WOTES:	Package (Y N Package (Y N Package (Y N N Package (Y N N N N N N N N N N N N N N N N N N	COC Record Was:  1 Present Upon-Receipt of Samples (Y N Discrepancies Between	Sample Labels and COC Record? Y N NOTES: 7-115
WESTON Analytics Use only  Custody Transfer Record/Lab Work Request	Wolume York / /	CINT. C.S.T.K.LILL F.F. ANALYSES	Collected Child With Children Collected Children X X X X X X X X X X X X X X X X X X X	2 sadvogative w x x x x x x x x x x x x x x x x x x	2 Sadioactive	so de certaine	W - Weiter DS - Drum Solids X - Other	 10-17	121-21-001/A-12/88 540094007 Less 4/101 200 102 //

EG&G ER Program Rocky Flats Plant

# ER PROGRAM DATA ASSESSMENT SUMMARY REPORT FORM

Bate	ch No. <u>8910L081</u>		Site _	Area 6 - Sola	r Ponds	·
Lab	oratory Roy F. Weston - Lionville		No. of	Samples/Ma	atrix <u>6/Wate</u>	<u> </u>
SOV	<b>X</b> # <u>7/87</u>		Review	wer Org. <u>Te</u>	chLaw, Inc.	
	nple Numbers <u>SW094007 (total), SW095007 (soluble), SW093007 (soluble</u>		al), SW09300	07 (total) SW	094007 (solub	le),
		Data Assess	ment Summa	ary		
		ICP	AA	Hg	CN	Comments
1.	Holding Times		<u>v</u>	<u>v</u>	<u>v</u>	
2.	Calibrations	A	<u>v</u>		v	Action Items 1-4
3.	Blanks	A	A		<u>v</u>	Action Items 5-10
4.	ICP Interference Check Sample	A	N/A	N/A	N/A	Action Items 11-13
5.	Lab Control Sample Results		<u>v</u>		v	
6.	Duplicate Sample Results	<u></u>	<u>V</u>	<u>v</u>	<u>v</u>	
7.	Matrix Spike Sample Results	v	A	<u>v</u>		Action Items 14-17
8.	Method of Standard Addition	N/A	<u>v</u>	N/A	N/A	
9.	Serial Dilution	_A	N/A	N/A	N/A	Action Item 18
10.	Sample Verification	v		v	<u>v</u>	
11.	Other QC	v			v	
12.	Overall Assessment	A	_A		V	Data valid, or acceptable with qualifications
	<ul> <li>V = Data had no problems.</li> <li>A = Data acceptable but qualified due to problems.</li> <li>R = Data rejected.</li> <li>X = Problems, but do not affect data.</li> </ul>				N/A = Not appl	icable.
Dat	ta Quality: Data contained in this batch w	vere reviewed ar	nd found to be v	alid, or acceptat	ole with qualifica	tions. Acceptable,
qual	lified data may be used provided that individu	ual values impac	cted by the "Act	ion Items" listed	i below are appro	priately flagged.
(Ref	fer to attached Results Summary Tables)					

Action Items: 1) All non-detect Molybdenum values are estimated and undetected (UJ) because the CRI
recovery criteria were not met.
2) The Zinc value for SW093007 (soluble) is estimated (J) because the CRI recovery criteria were not met.
3) The non-detect Lithium aqueous values for SW093007 (total and soluble) are estimated and undetected (UI)
because the CRI recovery criteria were not met.
4) All non-detect Tin values except SW093007 (total) are estimated and undetected (UJ) and the value for
SW093007 (total) is estimated (J) because the CRI recovery criteria were not met.
5) All Aluminum values are estimated and undetected (UJ) because Aluminum values >IDL were found in the
blanks.
6) The Cobalt values for SW094007 (total) and SW093007 (soluble) are estimated and undetected (UJ)
because Cobalt values >IDL were found in the blanks.
7) All Copper values are estimated and undetected (UJ) because Copper values >IDL were found in the
blanks.
8) The Potassium value for SW093007 (total and soluble) is estimated and undetected (UJ) because Potassium
values >IDL were found in the blanks.
9) The Antimony values for SW094007 (total and soluble), SW095007 (total), and SW093007 (total) are
estimated and undetected (UJ) because Antimony values >IDL were found in the blanks.
10) The Thallium value for SW095007 (soluble) is rejected (R) because of negative bias indicated in the
blanks.
11) The Silver values for SW094007 (total and soluble) and SW095007 (total and soluble) are rejected (R)
because the interference check sample result was outside control limits.
12) The Chromium values for SW094007 (total and soluble) and SW095007 (total and soluble) are rejected (R)
because the interference check sample result was outside control limits.
13) The Manganese, Beryllium, and Zinc values for SW094007 (total and soluble) and SW095007 (total and
soluble) are estimated (J) because the interference check sample result was outside control limits.

Action Items: (cont) 14) All Lead values except SW093007 (soluble) are estimated and undetected (UJ)
and the value for SW093007 (soluble) is estimated (J) because the pre-digestion matrix spike recovery criteria were
not met.
15) All Selenium values are estimated (J) because the pre-digestion matrix spike recovery criteria were not
met.
16) All Thallium values except SW093007 (total) are estimated and undetected (UJ) because the pre-digestion
matrix spike recovery criteria were not met.
17) The non-detect Arsenic values for SW094007 (soluble), SW095007 (soluble), and SW093007 (soluble) are
estimated and undetected (UJ) because the post-digestion matrix spike recovery criteria were not met.
18) All Calcium and Magnesium values are estimated (J) because the ICP serial dilution recovery criteria were
not met.
Comments: None
Comments. None
Note: Data Summary Tables are attached.
Reviewer Signature Date

SITE NAME: Site Background Characterization CLP WATER INORGANIC ANALYSIS: Low Water

Page 1 of 1

ANALYTICAL RESULTS (ug/L)

Sample Location							Н		Н					П
Sample Number		SW093007		SW093007		SW094007	S	SW094007	S	SW095007	SW095007	2000		
Sample Date		10/10/89		10/10/89		10/10/89	٥	10/10/89	2	10/10/89	10/10/89	66		
Remarks		Total		Soluble	Ħ	Total	S	Solubie	10	Total	Soluble			
Inorganic Analyte	DL Ug/L		g		g	۵	g	۵	g	g		g		
Aluminum Al	82	136 UJ	4	164 UJ	A 2	267 UJ	A 26	1 m 092	A 27	277 UJ A	293 UJ	٧		
	8	26.2 UJ	4	22.0 U	>	42.0 UJ	A 29	28.4 UJ /	A 39	39.9 UJ A	22.0 U	^		
Arsenic As	5	2.0 ∪	>	2.0 UJ	A	2.0 U	۷ 2.	2.0 UJ /	A 2.0	2.0 U V	2.0 W	4		
Barlum Ba	500	137	^	134	7	177	\ 	170	V 176	y .	1,1	>		
Beryllum Be	20	1.0 U	>	1.0 U	٧ ا	1.3 J	A 1.	1.3.J	A 1.:	1.3 J A	1.3 J	٧		
Cadmlum Cd	5	3.0 U	^	3.0 U	7	3.0 U	V 3.	3.0 ∪ \	3.0	3.0 U	3.0 ∪	>		
Calclum Ca	9009	112000 J	٧	109000 J	A	389000 J	A 3.	373000 J /	A 39	392000 J A	396000	A		
Cestum Ce	1000	2500 U	^	2500 U	۸ ک	2500 U	٧ ک	2500 U \	V 25	2500 U V	2500 U	>		
Chromlum Cr	5	2.0 U	>	2.0 U	۷ ا	2.0 U	R 2.	2.0 U	В 2.0	2.0 U R	2.0 U	н		
		4.0 U	^	4.4 UJ	4	5.1 UJ /	A 4.	4.0 U	۷ 4.(	4.0 U V	4.0 ∪	>		
Copper	52	11.5 UJ	٧	16.1 UJ	A	24.3 UJ	A 20	20.7 UJ /	A 25	25.1 UJ A	23.2 UJ	٧		
Iron F8	Ē	551	>	1220	۸	104	72		V 115	5 V	108	>		
Lead Pb	40	3.0 UJ	٧	85.6 J	A	3.0 UJ	A 3.	3.0 UJ	A 3.0	3.0 UJ A	3.0 UJ	4		
Lithium	0 <u>0</u>	100 UJ	٧	100 UJ	A	457	V 434		۷ 461	>	464	>		
Magnesium Mg	٦	30400 J	٧	29200 J	A	106000 J	4	102000 J	A 10	106000 J A	107000	4 T		
Manganese Mn		870		847	>	16.1 J		14.8.)	A 16	18.5 J A	16.1 J	∢		
Mercury Hg	0.2	0.20 U	>	0.20 U	>	0.20 U	о >	0.20 U	°; >	0.20 U	0.20 U	>		٦
Molybdenum Mo	200	100 UJ	٧	100 UJ	4	100 UJ	4 4	100 UJ /	4 4	100 UJ A	100 LU	٧		
Nickel Ni	40	7.0 U	>	7.0 U	>	9.3	<u>'</u>	7.0 U	우 >	10.3 V	7.0 U	>		٦
Potassium K	2000	3640 UJ	4	4300 UJ	<u>«</u>	91000	8 >	00298	9	91300 V	91200	>		
Selenium Se	ю	37.4.3	<	62.2 J	<u>-</u>	16.4 J	4	15.8.1	A 15	15.0 J A	923	V		
Silver Ag	0	3.0 U	>	3.0 U	>	3.0 U	F.	3.0 U	В.	3.0 U R	3.0 U	Œ		
Sodium Na	2000	61900	>	59600	>	000699	>	571000 \	-S >	V 000065	578000	>		
Strontium Sr	200	779	>	745	>	3120	× ×	3000	3	3120 V	3140	>		
Thailium TI	10	4.0 (J)	٧	4.0 UJ	A	40.0 UJ	4	4.0 UU /	A 4.	4.0 UJ A	4.0 U	Œ		
Tin Sn	88	100 UJ	4	100 UJ	4	136 J	4	114.3	A 12	128 J A	131	4		
Vanadlum V	8	5.0 U	>	5.0 U	>	5.0 U	> 2	5.0 U	);   2	5.0 U V	5.0 U	>		
Zinc Zn	8	65.6	>	54.3 J	<b>₹</b>	18.7 J	الا لا	23.4.1	N A	23.4.J A	24.8.)	4	-	
Cyanide	5	R		R.N	ᅴ	10.0 U	¥ ≥	н	뤼	10.0 U	ξ			7

E Estimated by the Laboratory
U indicates the compound was not detected above the instrument Quantitation Limit
J Quantitation is approximate due to firmitations identified during the quality control review
DL Detection Limit in Micrograms per Liker (ug/L)
NR Not reported

DQ Data Qualifier V Valid

A Acceptable with qualifications Rejected

L0811/eg04]

Lot 161089001 How STOCKTON MOTHULL

Custody Transfer Record/Lab Work Request

WISTON	WESTON Application	Use Only	Samples Were: (Shipped or Hand-	Delivered	NOTES:	NOTES:	3 Received Broken/	Leaking (Improperly	Sealed)	NOTES:	4 Roperly Preserved	**OTES:	5 Received Within	Thoding Times	₩OTES:		Present on Outer	- Artic	3 Present on Sample	•		COC Record Was:	of Samples (Y)	Incies E Labels	Record? Y (N NOTES:		7-115
el necord/Lab Work Request	1   6   6   5   5   5   1	er / 468 2148 445 / 1/45 1/1/49	7-17-17-17-	WON'S NOW JAN THURS HAR 3	7	S AND ON STATE OF SO MICES	海岛	2 X X X X X X	x x x x x x x x x x x x x x x x x x x	* × × ×	H		**			Kender 1020/89 Sanoko	A 4. 4 Tape NO Redinace			Scie West - Filtered tol metals Mo. Sr. Call a	unfiltered tol metals Mo, Sr. Ca, Li	Item/Reason Relinguished by Received by Date Time				20 000 MIND 000 EC	17 20 00 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1
ACAN OCTOR	Refrigerator#		[:	DO COLONIA DIA 1/10/10/10	ct/Phone	WA Use Only Client ID/Description Matrix Date	541094007 30 dination 14/ 1/2	X 18/0/089007	X 540 95007 radioaction	CCS 500 93007 SCOMBOOTHIS W	TO 100 P VS	ach Swaggood sadinacture	1			2	- L.20 CCP		X - Other		with the control of t	Time Time Time Time Time Time Time Time		1x (11)	13 mm 13 mm 1 1 1 1 1 1 1 1 1 1 1 1 1 1	FW 21-21-001/A-12/88	5/2001S

#### WESTON ANALYTICS TOTAL RADIOCHEMISTRY DATA SUMMARY REPORT

		Clien	t: ROCKWO	ELL (ROCKY FLATS)	)		Pa	зge: 3
Sample Information  RFW Batch ID: Customer ID: Collection Date: Matrix:	<b>%910%05</b> 9003-798-1 603F0689108900 10/05/89 Water	91		<b>891056787<sup>C</sup></b> 9003-798-2 604861089004 10/06/89 Water	")	8910108\ 9003-803-5 \$N094007 10/10/89 Nater		
Radio Chemistry  Gross Alpha Gross Beta Uranium 233, 234 Uranium 235 Uranium 238 Strontium 89, 90 Plutonium 239, 240 Americium 241 Cesium 137 Tritium Radium 226 Radimum 228	1.3 ± 0.7 3.3 ± 1.7 0.39 ± 0.39 0.10 ± 0.19 0.10 ± 0.19 0.11 ± 0.47 0.014 ± 0.008 0.006 ± 0.012 0.12 ± 0.44 40 ± 300		2.5 0 0 0.76 0 0.71	6.0 ± 2.0 8.7 ± 2.2 3.05 ± 1.09 0.09 ± 0.32 1.85 ± 0.87 0.68 ± 0.42 0.007 ± 0.008 0.012 ± 0.014 -0.18 ± 0.38 290 ± 280	pci/1 2.5 pci/1 2.8 pci/1 0.43 pci/1 0.43 pci/1 0.61 pci/1 0.010 pci/1 0.65 pci/1 430	146.1 ± 22.7 134.3 ± 17.7 84.9 ± 7.1 2.38 ± 1.26 49.7 ± 5.5 0.49 ± 0.44 0.017 ± 0.011 0.020 ± 0.010 0.11 ± 0.46 3430 ± 450	pci/l pci/l pci/l pci/l pci/l pci/l pci/l pci/l	17.9 0.69 0.69 0.69 0.67 0
	Province and an artist and an experience of the artist and artists and artists and artists and artists and art	Client	: ROCKWE	L <b>L (ROCKY</b> FLATS)				: sections
Sample Information  RFW Batch 10: Customer 10: Collection Date: Hatrix:	9003-803-6 SW095007 10/10/89 Water			9003-803-7 \$4093007 -10/10/89 Water				
Radio Chemistry  Gross Alpha. Gross Beta. Uranium 233, 234 Uranium 235. Uranium 238. Strontium 89, 90. Plutonium 239, 240. Americium 241. Cesium 137. Iritium. Radium 226. Radium 228.	120.4 ± 21.2 141.1 ± 18.8 70.2 ± 4.9 2.83 ± 1.02 45.3 ± 4.0 0.26 ± 0.42 0.010 ± 0.008 0.07 ± 0.42 2730 ± 400	pci/l pci/l pci/l pci/l pci/l pci/l pci/l pci/l	17.7 20.2 0.40 0.40 0.65 0.009 0.68 500	12.3 ± 2.7 13.4 ± 2.4 4.22 ± 1.17 0.08 ± 0.27 5.31 ± 1.31 0.77 ± 0.44 0.007 ± 0.005 0.010 ± 0.008 0.07 ± 0.41 -40 ± 260	pci/1 2.5 pci/1 2.7 pci/1 0.36 pci/1 0.36 pci/1 0.63 pci/1 0.63 pci/1 0			•